

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120215\
 Data File : BM003359.D
 Acq On : 02 Dec 2015 16:57
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM120215

Quant Time: Dec 02 17:37:21 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM120215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 02 17:30:44 2015
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	68	0.00
2	1,4-Dioxane	0.448	0.416	7.1	71	0.00
3	n-Nitrosodimethylamine	0.461	0.440	4.6	67	0.00
4 S	2-Fluorophenol	0.997	0.906	9.1	69	0.00
5 S	Phenol-d6	1.227	1.128	8.1	66	0.00
6	bis(2-Chloroethyl)ether	1.193	1.138	4.6	67	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	66	0.00
8 S	Nitrobenzene-d5	0.233	0.215	7.7	66	0.00
9	Nitrobenzene	0.252	0.232	7.9	64	0.00
10	Naphthalene	1.077	1.020	5.3	68	0.00
11	Hexachlorobutadiene	0.164	0.162	1.2	70	0.00
12	2-Methylnaphthalene	0.738	0.696	5.7	66	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	64	0.00
14 S	2,4,6-Tribromophenol	0.104	0.097	6.7	59	0.00
15 S	2-Fluorobiphenyl	1.505	1.388	7.8	66	0.00
16	Acenaphthylene	1.976	1.853	6.2	62	0.00
17	Acenaphthene	1.321	1.184	10.4	65	0.00
18	Fluorene	1.490	1.352	9.3	62	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	64	0.00
20	4-Bromophenyl-phenylether	0.144	0.123	14.6	63	0.00
21	Hexachlorobenzene	0.173	0.150	13.3	64	0.00
22	Pentachlorophenol	0.055	0.041	25.5#	52	0.00
23	Phenanthrene	0.964	0.865	10.3	64	0.00
24	Anthracene	1.016	0.911	10.3	58	0.00
25	Fluoranthene	0.999	0.902	9.7	57	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	51	0.00
27	Benzydine	0.409	0.366	10.5	52	0.02
28	Pvrene	1.812	1.774	2.1	55	0.00
29 S	Terphenyl-d14	0.811	0.829	-2.2	55	0.00
30	Benzo(a)anthracene	1.350	1.220	9.6	48#	0.00
31	3,3'-Dichlorobenzidine	0.349	0.294	15.8	45#	0.00
32	Chrysene	1.563	1.540	1.5	53	0.00
33	Bis(2-ethylhexyl)phthalate	0.684	0.561	18.0	42#	0.00
34	Indeno(1,2,3-cd)pyrene	1.163	1.042	10.4	59	0.01
35 I	Perylene-d12	1.000	1.000	0.0	52	0.00
36	Benzo(b)fluoranthene	1.522	1.413	7.2	51	0.00
37	Benzo(k)fluoranthene	1.366	1.254	8.2	50	0.00
38 C	Benzo(a)pyrene	1.249	1.102	11.8	50#	0.00
39	Dibenzo(a,h)anthracene	1.136	1.054	7.2	61	0.00
40	Benzo(a,h,i)perylene	1.199	1.096	8.6	61	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

(#) = Out of Range				
SPCC's out = 0 CCC's out = 0				